

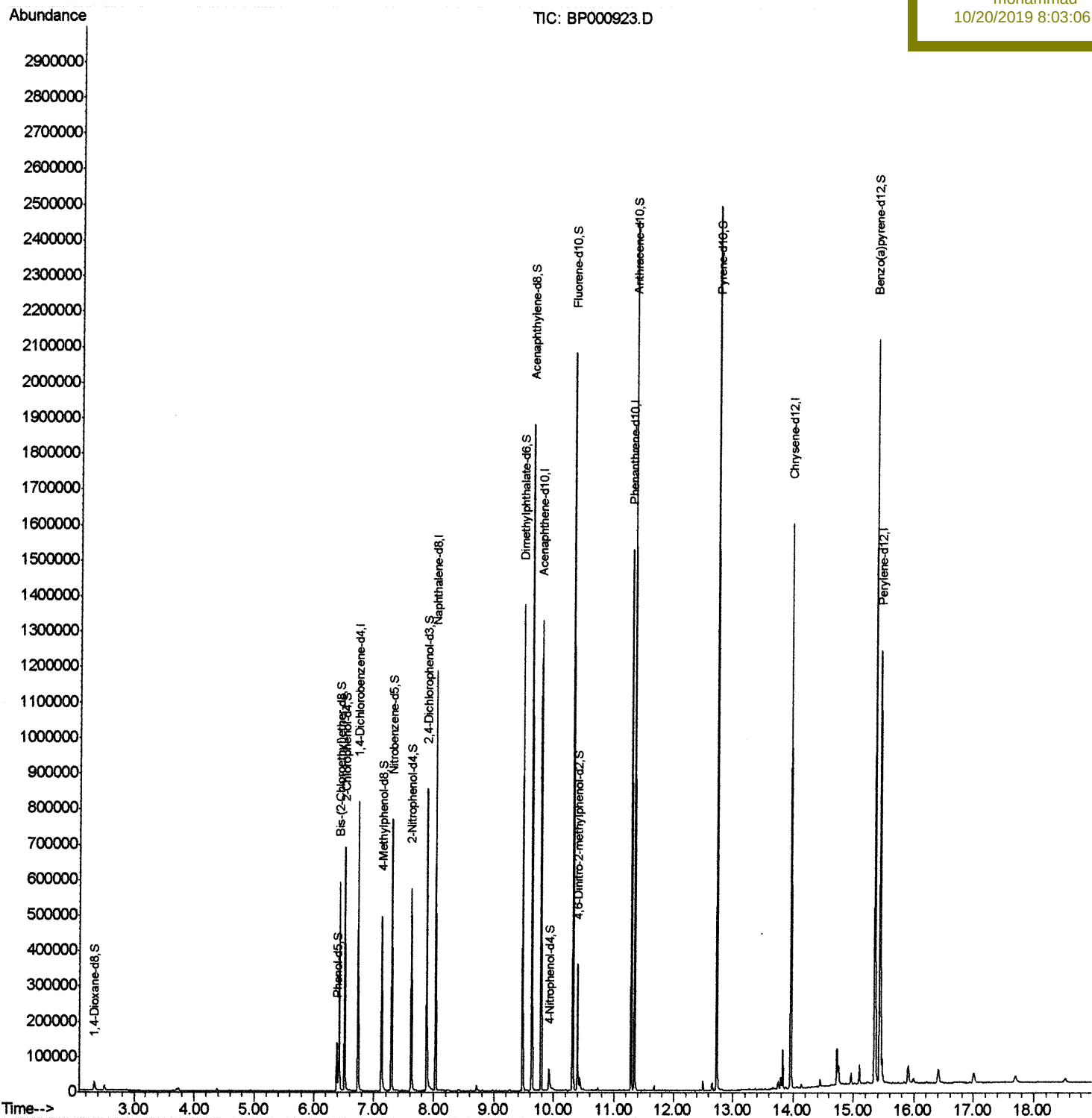
Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP101919\
Data File : BP000923.D
Acq On : 19 Oct 2019 18:39
Operator : JU
Sample : K5395-19
Misc :
ALS Vial : 21 Sample Multiplier: 1

Quant Time: Oct 19 22:59:24 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SOM-EPA-BP100919MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Sat Oct 19 22:14:11 2019
Response via : Initial Calibration

Instrument :
BNA_P
ClientSampleId :
C0052

Manual Integrations
APPROVED

mohammad
10/20/2019 8:03:06 PM



Quantitation Report (Qedit)

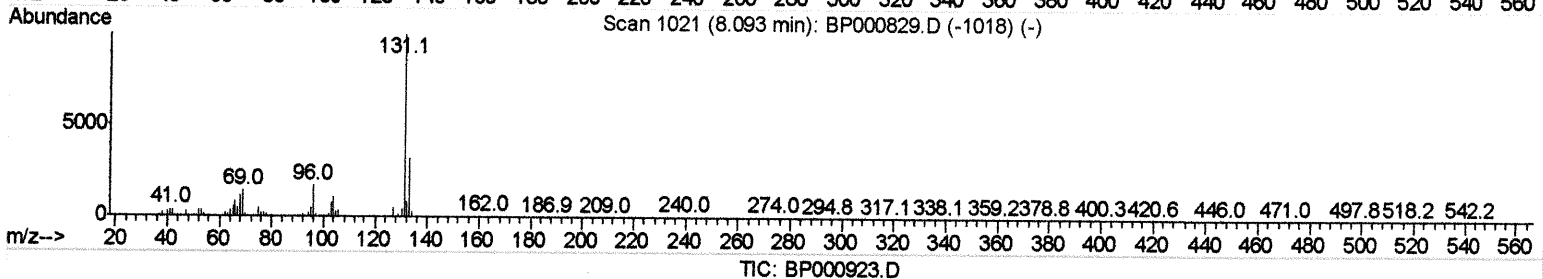
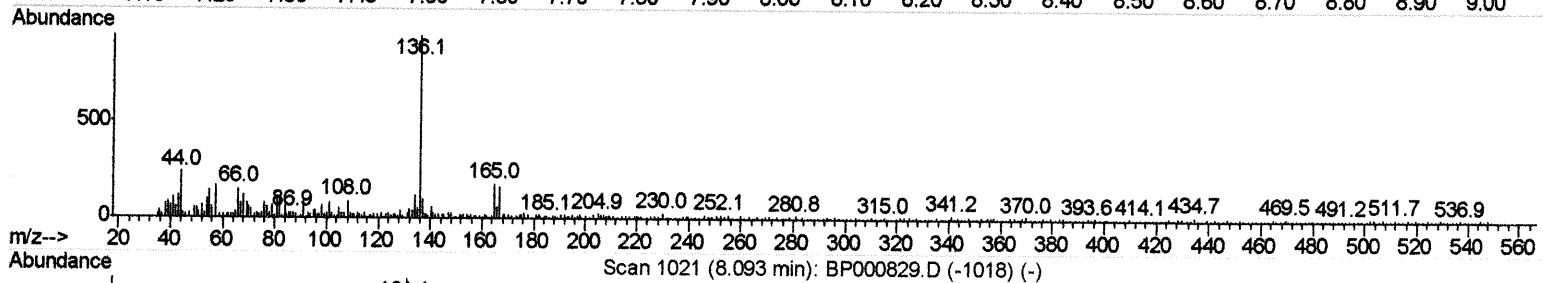
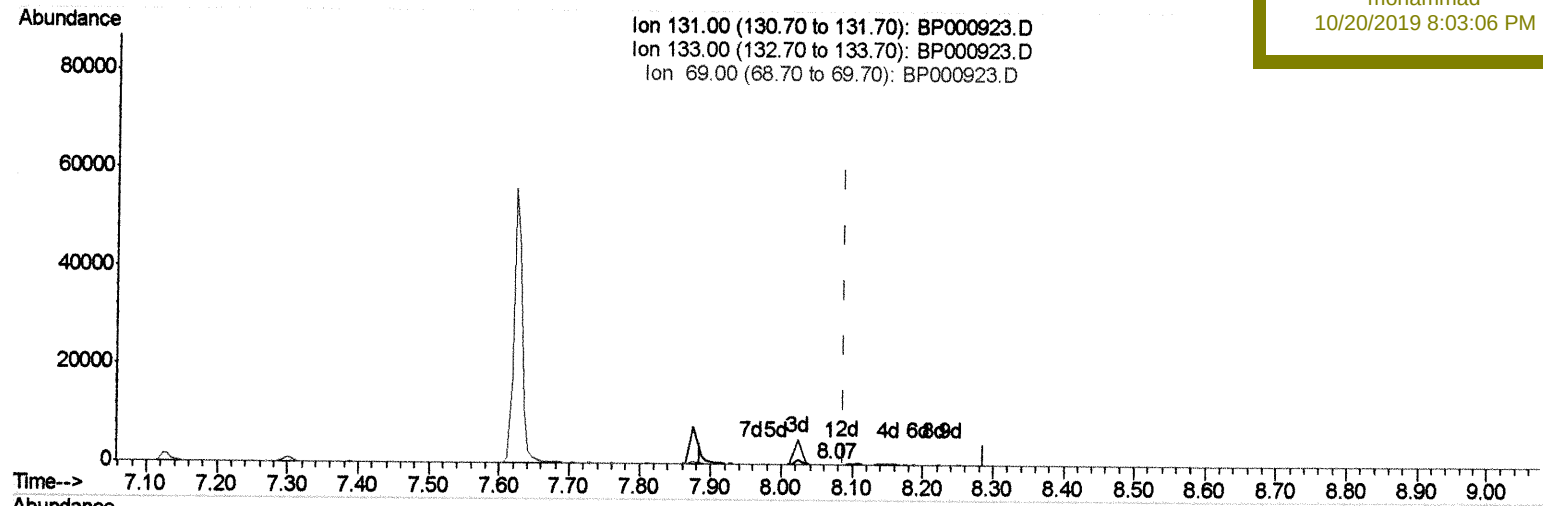
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Quant Time: Oct 19 22:58:29 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SOM-EPA-BP100919MA.M
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(29) 4-Chloroaniline-d4 (S)

8.069min (-0.018) 0.00ng/ul

response 6

Ion	Exp%	Act%
131.00	100	100
133.00	32.70	148.15#
69.00	13.70	277.78#
0.00	0.00	0.00

Quantitation Report (Qedit)

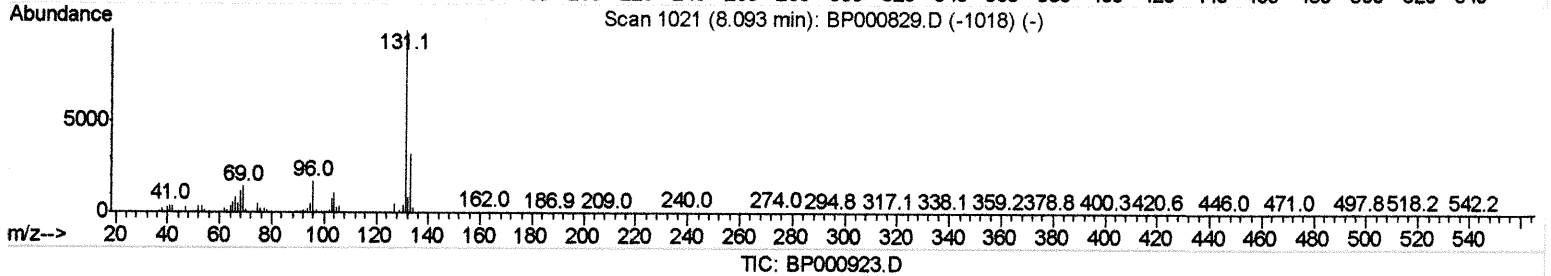
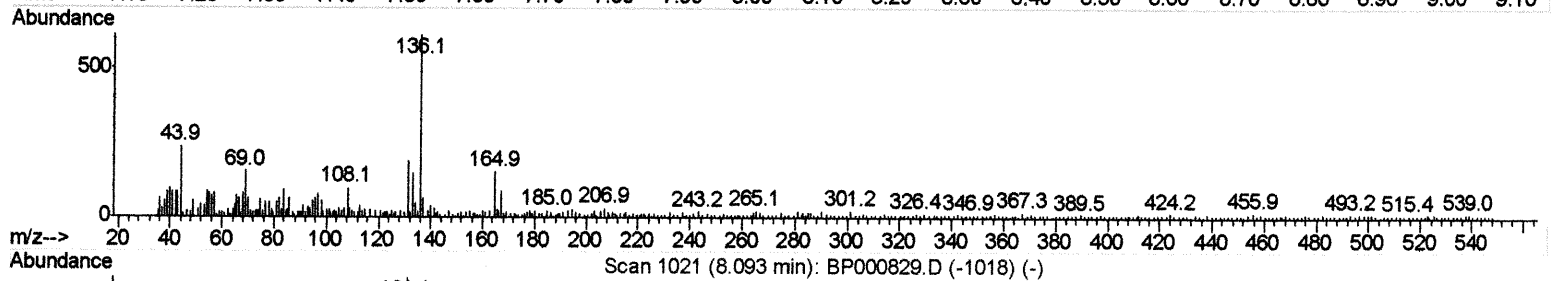
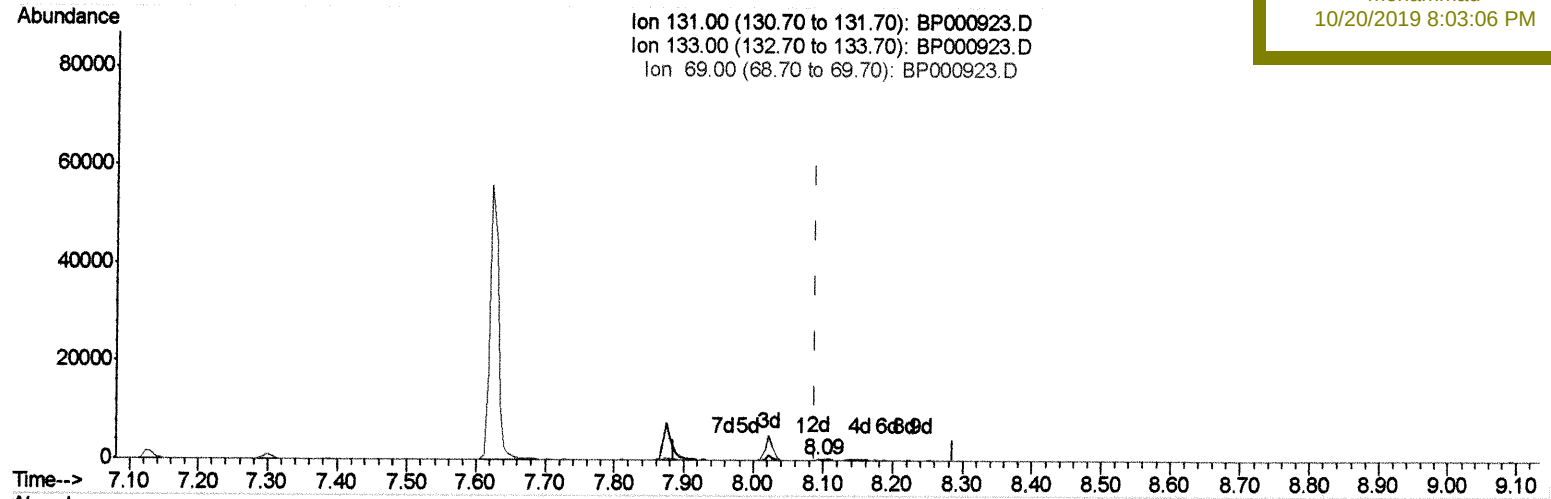
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Manual Integrations
 APPROVED

mohammad
 10/20/2019 8:03:06 PM



(29) 4-Chloroaniline-d4 (S)

8.093min (+0.006) 0.03ng/ul m

response 277

Ion	Exp%	Act%
131.00	100	100
133.00	32.70	77.13#
69.00	13.70	82.45#
0.00	0.00	0.00

JUC-11/19

Data Path : Z:\SVOASRV\HPCHEM1\BNA_P\DATA\BP101919\
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 Acq On : 19 Oct 2019 18:39
 Operator : JU
 Sample : K5395-19
 Misc :
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 BNA_P
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Manual Integrations
 APPROVED

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev (Min)
1) 1,4-Dichlorobenzene-d4	6.74	152	122542	20.00	ng/ul	0.00
18) Naphthalene-d8	8.02	136	461323	20.00	ng/ul	0.00
36) Acenaphthene-d10	9.79	164	262327	20.00	ng/ul	0.00
62) Phenanthrene-d10	11.28	188	512557	20.00	ng/ul	0.00
78) Chrysene-d12	13.96	240	496043	20.00	ng/ul	0.00
86) Perylene-d12	15.43	264	562337	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	2.34	96	15287	5.17	ng/uL	0.00
5) Phenol-d5	6.39	99	59706	6.32	ng/ul	0.00
7) Bis-(2-Chloroethyl) ether-d	6.42	67	123879	26.21	ng/ul	0.00
9) 2-Chlorophenol-d4	6.51	132	187603	23.56	ng/ul	0.00
13) 4-Methylphenol-d8	7.13	113	116900	16.19	ng/ul	0.00
19) Nitrobenzene-d5	7.30	128	105398	30.06	ng/ul	0.00
22) 2-Nitrophenol-d4	7.62	143	117972	31.59	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	7.88	165	196143	27.44	ng/ul	0.00
29) 4-Chloroaniline-d4	8.09	131	277m	0.03	ng/ul	0.00
44) Dimethylphthalate-d6	9.48	166	614514	33.46	ng/ul	0.00
47) Acenaphthylene-d8	9.63	160	701643	31.06	ng/ul	0.00
52) 4-Nitrophenol-d4	9.92	143	17274	6.11	ng/ul	0.00
58) Fluorene-d10	10.30	176	520184	33.11	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	10.39	200	58654	24.03	ng/ul	0.00
71) Anthracene-d10	11.34	188	808013	34.03	ng/ul	0.00
79) Pyrene-d10	12.72	212	884324	31.53	ng/ul	0.00
90) Benzo(a)pyrene-d12	15.35	264	960506	34.60	ng/ul	0.00

2019/10/20

Target Compounds Ovalue

(#) = qualifier out of range (m) = manual integration (+) = signals summed